original publications.

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FILE COVERS 1907 - 13 Apr 2007 VOL 146 ISS 17 FILE LAST UPDATED: 12 Apr 2007 (20070412/ED)

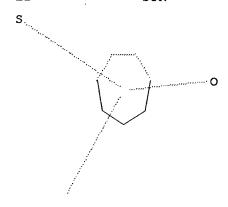
Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

http://www.cas.org/infopolicy.html

Uploading C:\Program Files\Stnexp\Queries\781.str

L1 STRUCTURE UPLOADED

=> d L1 HAS NO ANSWERS STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1 full REG1stRY INITIATED

Substance data SEARCH and crossover from CAS REGISTRY in progress... Use DISPLAY HITSTR (or FHITSTR) to directly view retrieved structures.

FULL SEARCH INITIATED 15:23:13 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 6914556 TO ITERATE

6.5% PROCESSED 450880 ITERATIONS

20 ANSWERS

11.7% PROCESSED 810862 ITERATIONS

30 ANSWERS

14.5% PROCESSED 1000000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

32 ANSWERS

SEARCH TIME: 00.00.44

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**

BATCH **INCOMPLETE**

PROJECTED ITERATIONS:

6914556 TO 6914556

PROJECTED ANSWERS:

177 TO 265

L2

32 SEA SSS FUL L1

L3

8 L2

=> s 13 and PY<2002

21882609 PY<2002

T.4

0 L3 AND PY<2002

=> d 13 1-8 ibib abs hitstr

L3 ANSWER 1 OF 8 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER:

2007:227224 CAPLUS

DOCUMENT NUMBER:

146:268408

TITLE:

Phenylsulfonylcarbamate derivatives as herbicide

safeners

INVENTOR(S):

Furuse, Katsumi; Takahashi, Satoru; Ohno, Shuji;

Ogawa, Yasunori; Mitsunari, Takashi

PATENT ASSIGNEE(S):

Kumiai Chemical Industry Co., Ltd., Japan

SOURCE:

PCT Int. Appl., 78pp. CODEN: PIXXD2

DOCUMENT TYPE:

Patent

1

Ι

LANGUAGE:

GI

Japanese

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PAT	PATENT NO.					D	DATE		APPLICATION NO.						DATE			
₩O	2007023764				A1 20070301			WO 2006-JP316316						20060821				
	W:	ΑE,	AG,	AL,	AM,	AT,	ΑU,	ΑZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH.	
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD.	
•		GE,	GH,	GM,	HN,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KM,	KN,	KP,	
		KR,	ΚZ,	LA,	LC,	LK,	LR,	LS,	LT,	LU,	LV,	LY,	MA,	MD,	MG,	MK,	MN.	
		MW,	MX,	MY,	ΜZ,	NΑ,	NG,	NI,	NO,	NZ,	OM,	PG,	PH,	PL,	PT,	RO,	RS,	
		RU,	SC,	SD,	SE,	SG,	SK,	SL,	SM,	SV,	SY,	ТJ,	TM,	TN,	TR,	TT,	TZ,	
		UA,	ŪG,	US,	UZ,	VC,	VN,	ZA,	ZM,	ZW					-		-	
	RW:	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,	FI,	FR,	GB,	GR,	HU,	IE,	
		IS,	IT,	LT,	LU,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR,	BF,	ВJ,	
		CF,	CG,	CI,	CM,	GΑ,	GN,	GQ,	GW,	ML,	MR,	NE,	SN,	TD,	TG,	BW,	GH,	
		GM,	ΚE,	LS,	MW,	MZ,	NA,	SD,	SL,	SZ,	TZ;	ŬĠ,	ZM,	ZW,	AM,	AZ,	BY,	
				MD,											·	•	•	
PRIORITY	APP	LN.	INFO	.:					JP 2005-245544					A 20050826				

$$y_{n} = \begin{cases} co - x_{R}1 \\ so_{2} - N - co - z_{R}2 \\ l \\ R3 \end{cases}$$

AB Phenylsulfonylcarbamate derivs. (I, wherein R1, R2, R3 = H, alkyl, etc.; Y = halo, NO2, etc.; n = 0-4 integer; X, Z = 0, S) or salts thereof decrease

the harmful effect of herbicides against cultivated plants without loss of effectiveness. Thus, when benzobicyclon (40 g/10 are) was applied 5 days after transplanting rice in a pot experiment, growth inhibition was 20% at 29 days after transplanting, whereas when I (R1, R3 = H, R2 = 4-chlorobenzyl, X = 0, n = 0) was applied at 240 g/10 are on the day after transplanting with the same benzobicyclon treatment, the growth inhibition with only 8%. 927411-99-2

RL: AGR (Agricultural use); BSU (Biological study, unclassified); BIOL (Biological study); USES (Uses)

(benzobicyclon + Bensulfuron-Me + compound III-1; safened herbicide composition)

RN 927411-99-2 CAPLUS

CN Benzoic acid, 2-[[[[(4,6-dimethoxy-2-pyrimidinyl)amino]carbonyl]amino]sul fonyl]methyl]-, methyl ester, mixt. with 3-[2-chloro-4-(methylsulfonyl)benzoyl]-4-(phenylthio)bicyclo[3.2.1]oct-3-en-2-one and methyl 2-[[(ethoxycarbonyl)amino]sulfonyl]benzoate (CA INDEX NAME)

CM 1

ΙT

CRN 156963-66-5 CMF C22 H19 Cl O4 S2

CM 2

CRN 83404-84-6 CMF C11 H13 N O6 S

CM 3

CRN 83055-99-6 CMF C16 H18 N4 O7 S

IT 927411-95-8

RL: AGR (Agricultural use); BSU (Biological study, unclassified); BIOL (Biological study); USES (Uses)

(benzobicyclon + cafenstrole + compound I-36; safened herbicide composition)

RN 927411-95-8 CAPLUS

CN Benzoic acid, 2-[[[(4-chlorophenyl)methoxy]carbonyl]amino]sulfonyl]-,
mixt. with 3-[2-chloro-4-(methylsulfonyl)benzoyl]-4(phenylthio)bicyclo[3.2.1]oct-3-en-2-one and N,N-diethyl-3-[(2,4,6trimethylphenyl)sulfonyl]-1H-1,2,4-triazole-1-carboxamide (CA INDEX NAME)

CM 1

CRN 808197-84-4 CMF C15 H12 C1 N O6 S

CM 2

CRN 156963-66-5 CMF C22 H19 Cl O4 S2

CM 3

CRN 125306-83-4 CMF C16 H22 N4 O3 S

$$\begin{array}{c|c} O & Me \\ \parallel & Me \\ Et_2N-C & N & \parallel \\ N & O & Me \\ \end{array}$$

IT 927411-91-4

RL: AGR (Agricultural use); BSU (Biological study, unclassified); BIOL (Biological study); USES (Uses)

(benzobicyclon + compound I-35; safened herbicide composition)

RN 927411-91-4 CAPLUS

CN Benzoic acid, 2-[[[[(2-chlorophenyl)methoxy]carbonyl]amino]sulfonyl]-, mixt. with 3-[2-chloro-4-(methylsulfonyl)benzoyl]-4- (phenylthio)bicyclo[3.2.1]oct-3-en-2-one (CA INDEX NAME)

CM 1

CRN 808197-83-3 CMF C15 H12 C1 N O6 S

CM 2

CRN 156963-66-5 CMF C22 H19 C1 O4 S2

IT 927411-88-9

RL: AGR (Agricultural use); BSU (Biological study, unclassified); BIOL (Biological study); USES (Uses)

(benzobicyclon + compound I-36; safened herbicide composition)

RN 927411-88-9 CAPLUS

CN Benzoic acid, 2-[[[[(4-chlorophenyl)methoxy]carbonyl]amino]sulfonyl]-, mixt. with 3-[2-chloro-4-(methylsulfonyl)benzoyl]-4- (phenylthio)bicyclo[3.2.1]oct-3-en-2-one (CA INDEX NAME)

CM 1

CRN 808197-84-4 CMF C15 H12 C1 N O6 S

CM2

CRN 156963-66-5 CMF C22 H19 C1 O4 S2

REFERENCE COUNT:

10 THERE ARE 10 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

CAPLUS COPYRIGHT 2007 ACS on STN ANSWER 2 OF 8

ACCESSION NUMBER:

2007:226871 CAPLUS

DOCUMENT NUMBER:

146:268407

TITLE: INVENTOR(S):

Benzoisothiazolinone dioxides as herbicide safeners

Furuse, Katsumi; Ueno, Ryohei; Asakura, Sohei; Yonekura, Norihisa; Mitsunari, Takashi

PATENT ASSIGNEE(S):

Kumiai Chemical Industry Co., Ltd., Japan

SOURCE:

PCT Int. Appl., 68pp.

CODEN: PIXXD2

DOCUMENT TYPE:

Patent

LANGUAGE:

GI

Japanese

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PAT	PATENT NO.					D	DATE		APPLICATION NO.						DATE				
WO	WO 2007023719				A1 20070301			WO 2006-JP316097						20060816					
	W:	ΑE,	AG,	AL,	AM,	ΑT,	AU,	ΑZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ.	CA.	CH.		
		CN,	co,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD.		
		GE,	GH,	GM,	HN,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KM,	KN.	KP.		
		KR,	KZ,	LА,	LC,	LK,	LR,	LS,	LT,	LU,	LV,	LY,	MA,	MD,	MG,	MK.	MN.		
		MW,	MX,	MY,	MZ,	NA,	NG,	NI,	NO,	ΝZ,	OM,	PG,	PH,	PL,	PT,	RO,	RS.		
		RU,	SC,	SD,	SE,	SG,	SK,	SL,	SM,	SY,	ТJ,	TM,	TN,	TR,	TT,	TZ,	UA,		
		ŪG,	US,	UZ,	VC,	VN,	ZA,	ZM,	ZW							•	•		
	RW:	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,	FI,	FR,	GB,	GR,	HU,	IE,		
		IS,	IT,	LT,	LU,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR,	BF,	BJ,		
		CF,	CG,	CI,	CM,	GA,	GN,	GQ,	GW,	ML,	MR,	NE,	SN,	TD,	TG,	BW,	GH,		
		GM,	ΚE,	LS,	MW,	MZ,	NA,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	AM,	AZ,	BY.		
		KG,	KZ,	MD,	RU,	TJ,	TM								·	•	•		
PRIORITY	APP	LN.	INFO	.:					JP 2005-239757					Z	A 20050822				

$$x_n$$
 x_n
 x_n

AB 1,2-Benzoisothiazolin-3-one-1,1-dioxide derivs. (I, wherein Y = 0, S; R1 = C1-16 alkyl, C2-6 alkenyl, etc.; X = halo, NO2, alkyl, etc.; n = 0-4 integer) or salts thereof are extremely favorable for reducing chemical injury to cultivated plants without reducing weed control by herbicides. Thus, in a pot experiment I (Y = 0, R1 = 4-chlorobenzyl, n = 0) was applied at 240 g/10 are on the day after transplanting (DAT) of rice, and benzobicyclon was applied at 20 g/10 are at 5 DAT. There was no inhibition of rice growth at 32 DAT, whereas rice growth inhibition was 6% when benzobicyclon was applied without the safener. In another experiment with benzobicyclon applied at 12.5 g/10 are, control of Scirpus juncoides was ≥90%, whether or not pots were pretreated with 240 g/10 are of the same I derivative

IT 927419-12-3 927419-15-6 927419-22-5

I

927419-26-9

RL: AGR (Agricultural use); BSU (Biological study, unclassified); BIOL (Biological study); USES (Uses)

(safened herbicide compns.)

RN 927419-12-3 CAPLUS

CN 1,2-Benzisothiazole-2(3H)-carboxylic acid, 3-oxo-, (4-chlorophenyl)methyl ester, 1,1-dioxide, mixt. with 3-[2-chloro-4-(methylsulfonyl)benzoyl]-4-(phenylthio)bicyclo[3.2.1]oct-3-en-2-one (CA INDEX NAME)

CM 1

CRN 863554-50-1 CMF C15 H10 C1 N O5 S

CM 2

CRN 156963-66-5 CMF C22 H19 C1 O4 S2

RN 927419-15-6 CAPLUS

CN 1,2-Benzisothiazole-2(3H)-carboxylic acid, 3-oxo-, (2-chlorophenyl)methyl ester, 1,1-dioxide, mixt. with 3-[2-chloro-4-(methylsulfonyl)benzoyl]-4-(phenylthio)bicyclo[3.2.1]oct-3-en-2-one (CA INDEX NAME)

CM 1

CRN 927419-03-2 CMF C15 H10 C1 N O5 S

CM 2

CRN 156963-66-5 CMF C22 H19 C1 O4 S2

RN 927419-22-5 CAPLUS

CN 1,2-Benzisothiazole-2(3H)-carboxylic acid, 3-oxo-, (4-chlorophenyl)methyl ester, 1,1-dioxide, mixt. with 3-[2-chloro-4-(methylsulfonyl)benzoyl]-4-(phenylthio)bicyclo[3.2.1]oct-3-en-2-one and N,N-diethyl-3-[(2,4,6-trimethylphenyl)sulfonyl]-1H-1,2,4-triazole-1-carboxamide (CA INDEX NAME)

CM 1

CRN 863554-50-1 CMF C15 H10 C1 N O5 S

CM 2

CRN 156963-66-5

CM 3

CRN 125306-83-4 CMF C16 H22 N4 O3 S

$$\mathsf{Et}_2\mathsf{N}-\mathsf{C} = \mathsf{N} = \mathsf{N} = \mathsf{N} = \mathsf{Me}$$

RN 927419-26-9 CAPLUS

CN 1,2-Benzisothiazole-2(3H)-carboxylic acid, 3-oxo-, (4-chlorophenyl)methyl ester, 1,1-dioxide, mixt. with 3-[2-chloro-4-(methylsulfonyl)benzoyl]-4-(phenylthio)bicyclo[3.2.1]oct-3-en-2-one and methyl 2-[[[[(4,6-dimethoxy-2-pyrimidinyl)amino]carbonyl]amino]sulfonyl]methyl]benzoate (CA INDEX NAME)

CM 1

CRN 863554-50-1 CMF C15 H10 C1 N O5 S

CM 2

CRN 156963-66-5 CMF C22 H19 C1 O4 S2

CM 3

CRN 83055-99-6 CMF C16 H18 N4 O7 S

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L3 ANSWER 3 OF 8 CAPLUS COPYRIGHT 2007 ACS on STN ACCESSION NUMBER: 2007:120952 CAPLUS

7

DOCUMENT NUMBER:

REFERENCE COUNT:

146:206022

TITLE:

Synthetic studies on the MARDi cascade:

stereoselective preparation of sulfonyl-substituted seven-membered rings. [Erratum to document cited in

THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS

CA146:100354]

AUTHOR(S):

Coquerel, Yoann; Bensa, David; Moret, Vincent;

Rodriguez, Jean

CORPORATE SOURCE:

UMR CNRS 6178, Centre Universitaire de St. Jerome,

Universite Paul Cezanne (Aix-Marseille III),

Marseille, 13397/20, Fr.

SOURCE: Synlett (2006), (19), 3368 CODEN: SYNLES; ISSN: 0936-5214

PUBLISHER:

Georg Thieme Verlag Journal

DOCUMENT TYPE:

English

LANGUAGE:

On page 2752, the chemical structure of cycloheptanol as compound (8) in Table ΑB 1 was incorrectly represented. The correct structure is given.

IT 917971-71-2P 917971-72-3P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (stereoselective synthesis of functionalized sulfonyl-substituted cycloheptanes via formal two-carbon ring expansion of 2-benzenesulfonyl cyclopentanones through a base-induced anionic domino three-component transformation (Erratum))

RN 917971-71-2 CAPLUS

Cycloheptanecarboxylic acid, 2-hydroxy-4-methyl-5-(phenylsulfonyl)-, CN methyl ester, (1R,2R,4R,5R)-rel- (CA INDEX NAME)

Relative stereochemistry.

RN 917971-72-3 CAPLUS

CN Cycloheptanecarboxylic acid, 2-hydroxy-3-methyl-5-(phenylsulfonyl)-, methyl ester, (1R,2R,3S,5R)-rel- (CA INDEX NAME)

Relative stereochemistry.

IT 917971-70-1P 917971-73-4P

RL: SPN (Synthetic preparation); PREP (Preparation)
(stereoselective synthesis of functionalized sulfonyl-substituted
cycloheptanes via formal two-carbon ring expansion of 2-benzenesulfonyl
cyclopentanones through a base-induced anionic domino three-component
transformation (Erratum))

RN 917971-70-1 CAPLUS

CN Cycloheptanecarboxylic acid, 2-hydroxy-5-(phenylsulfonyl)-, methyl ester, (1R,2R,5R)-rel- (CA INDEX NAME)

Relative stereochemistry.

RN 917971-73-4 CAPLUS

CN Cycloheptanecarboxylic acid, 3-butyl-2-hydroxy-5-(phenylsulfonyl)-, methyl ester, (1R,2R,3S,5R)-rel- (CA INDEX NAME)

Relative stereochemistry.

L3 ANSWER 4 OF 8 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER:

2006:1261562 CAPLUS

DOCUMENT NUMBER:

146:206186

TITLE:

Polycyclic oxonium ylides - Use of cyclic acetals as convenient scaffolds in the construction of fused

bicyclic compounds containing a medium ring

AUTHOR(S):

CORPORATE SOURCE:

Murphy, Graham K.; Marmsaeter, Fredrik P.; West, F. G. Department of Chemistry, Gunning-Lemieux Chemistry

Centre, University of Alberta, Edmonton, AB, T6G 2G2,

SOURCE:

Canadian Journal of Chemistry (2006), 84(10),

II

1470-1486

CODEN: CJCHAG; ISSN: 0008-4042

National Research Council of Canada

PUBLISHER:

DOCUMENT TYPE:

LANGUAGE:

GΙ

English

I

Cyclic mixed acetals and thioacetals I (R = H, MeO, 4-MeC6H4S; R1 = MeO, AB 4-MeC6H4S, H; R2 = H, EtO2C; X = CH2, CH2CH2) with pendant diazoketones undergo efficient rearrangement to ether-bridged cyclooctanoid and cycloheptanoid systems such as oxatricycles II (R = H, MeO, 4-MeC6H4S; R1 = MeO, 4-MeC6H4S, H; R2 = H, EtO2C; X = CH2, CH2CH2) upon treatment with copper bis(hexafluoroacetylacetonate). Other catalysts such as copper bis(trifluoroacetylacetonate), dirhodium tetraacetate, and dirhodium tetrakis(triphenylacetate) are significantly less effective in generating oxygen-bridged polycycles from I. A mechanism for the cyclocondensation is proposed; generation of oxonium ylides from I is followed by a [1,2]-shift to generate II. This work indicates that heteroatomsubstituted oxonium ylides can undergo Stevens [1,2]-shifts. The arylthio moiety of products derived from mixed thioacetals can either be reductively cleaved or can be used to cleave the bridging ether. IT 923054-48-2P

RL: BYP (Byproduct); PREP (Preparation) (byproduct in the stereoselective preparation of oxatricycles by ylide formation and stereoselective rearrangement of diazoketones containing cyclic mixed acetals and thioacetals)

RN 923054-48-2 CAPLUS

CN 1H-3a,7-Epoxyazulene-6-carboxylic acid, octahydro-6-[(4-methylphenyl)thio]-5-oxo-, ethyl ester, (3aR,6S,7R,8aS)-rel- (CA INDEX NAME)

Relative stereochemistry.

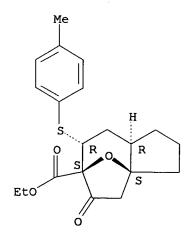
IT 923054-45-9P

RL: SPN (Synthetic preparation); PREP (Preparation) (stereoselective preparation of oxatricycles by ylide formation and stereoselective rearrangement of diazoketones containing cyclic mixed acetals and thioacetals)

RN 923054-45-9 CAPLUS

CN 6H-3a,6-Epoxyazulene-6-carboxylic acid, octahydro-7-[(4-methylphenyl)thio]-5-oxo-, ethyl ester, (3aR,6R,7S,8aS)-rel- (CA INDEX NAME)

Relative stereochemistry.



REFERENCE COUNT:

THERE ARE 44 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L3 ANSWER 5 OF 8 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER:

2006:1188387 CAPLUS

DOCUMENT NUMBER:

146:100354

TITLE:

Synthetic studies on the MARDi cascade:

stereoselective preparation of sulfonyl-substituted

seven-membered rings

AUTHOR(S):

Coquerel, Yoann; Bensa, David; Moret, Vincent;

Rodriguez, Jean

CORPORATE SOURCE:

UMR CNRS 6178, Centre Universitaire de St Jerome,

Universite Paul Cezanne (Aix-Marseille III),

Marseille, 13397/20, Fr.

SOURCE:

Synlett (2006), (17), 2751-2754

CODEN: SYNLES; ISSN: 0936-5214

PUBLISHER: Georg Thieme Verlag

DOCUMENT TYPE:

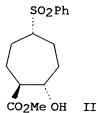
Journal

LANGUAGE:

English

GI

SO₂Ph



AB A stereoselective synthesis of functionalized sulfonyl-substituted cycloheptanes is described. The approach involves a formal two-carbon ring expansion of 2-benzenesulfonyl cyclopentanones through a base-induced anionic domino three-component transformation named the MARDi cascade (Michael Aldol Retro-Dieckmann). E.g., to a solution of β -keto sulfone I was added CH2:CHCHO and K2CO3 to give 62% cycloheptane II (dr 4:1). IT 917971-71-2P 917971-72-3P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (stereoselective synthesis of functionalized sulfonyl-substituted cycloheptanes via formal two-carbon ring expansion of 2-benzenesulfonyl cyclopentanones through a base-induced anionic domino three-component transformation)

RN 917971-71-2 CAPLUS

CN Cycloheptanecarboxylic acid, 2-hydroxy-4-methyl-5-(phenylsulfonyl)-, methyl ester, (1R,2R,4R,5R)-rel- (CA INDEX NAME)

Relative stereochemistry.

RN 917971-72-3 CAPLUS

CN Cycloheptanecarboxylic acid, 2-hydroxy-3-methyl-5-(phenylsulfonyl)-, methyl ester, (1R,2R,3S,5R)-rel- (CA INDEX NAME)

Relative stereochemistry.

IT 917971-70-1P 917971-73-4P

> RL: SPN (Synthetic preparation); PREP (Preparation) (stereoselective synthesis of functionalized sulfonyl-substituted cycloheptanes via formal two-carbon ring expansion of 2-benzenesulfonyl cyclopentanones through a base-induced anionic domino three-component transformation)

917971-70-1 CAPLUS RN

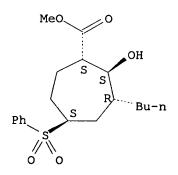
Cycloheptanecarboxylic acid, 2-hydroxy-5-(phenylsulfonyl)-, methyl ester, CN (1R, 2R, 5R) - rel - (CA INDEX NAME)

Relative stereochemistry.

RN 917971-73-4 CAPLUS

Cycloheptanecarboxylic acid, 3-butyl-2-hydroxy-5-(phenylsulfonyl)-, methyl CN ester, (1R, 2R, 3S, 5R) -rel- (CA INDEX NAME)

Relative stereochemistry.



REFERENCE COUNT: 19 THERE ARE 19 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L3 CAPLUS COPYRIGHT 2007 ACS on STN ANSWER 6 OF 8

ACCESSION NUMBER:

2006:744227 CAPLUS

TITLE:

Chemical components of essential oils from Liquidambar

orientalis Mill

AUTHOR(S):

Yao, Faye; Qiu, Qin; Cui, Zhaojie; Su, Demin Department of Chemistry, Shandong Institute of

CORPORATE SOURCE:

Education, Jinan, 250013, Peop. Rep. China

SOURCE:

Yaowu Fenxi Zazhi (2005), 25(7), 859-862

CODEN: YFZADL; ISSN: 0254-1793

Yaowu Fenxi Zazhi Bianji Weiyuanhui

PUBLISHER: DOCUMENT TYPE:

Journal

LANGUAGE: Chinese

To analyze the chemical constituents of volatile oil from the root of Liquidambar orientalis Mill, the volatile oils from roots of Liquidambar orientalis Mill was extracted by SFE CO2, and analyzed by gas chromatog.-mass spectrometry (GC-MS). Fifty compds. were identified, which represented 87% of the total GC peak area of the volatile oil. The present study provides scientific bases for the Liquidambar orientalis Mill exploitation in reason.

929903-86-6 IT

> RL: COS (Cosmetic use); BIOL (Biological study); USES (Uses) (chemical constituents of volatile oils of Liquidambar)

RN 929903-86-6 CAPLUS

CN INDEX NAME NOT YET ASSIGNED

OH Me Me PhS Me

CAPLUS COPYRIGHT 2007 ACS on STN ANSWER 7 OF 8

ACCESSION NUMBER:

2006:600049 CAPLUS

DOCUMENT NUMBER:

145:248986

TITLE:

Double Lawton SN2' Addition to Epoxyvinyl Sulfones:

Selective Construction of the Stereotetrads of

Aplyronine A

AUTHOR(S):

El-Awa, Ahmad; Fuchs, Philip

CORPORATE SOURCE:

Department of Chemistry, Purdue University, West

Lafayette, IN, 47907, USA

SOURCE:

Organic Letters (2006), 8(14), 2905-2908

CODEN: ORLEF7; ISSN: 1523-7060

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CASREACT 145:248986

GT

AB Enantiopure epoxyvinyl sulfones I (R = Me, R1 = H; R = H, R1 = Me) function as templates for the diastereoselective construction of the three stereotetrads of aplyronine A. Lawton SN2' addition of 3,5-dimethylpyrazole followed by its displacement in an alc.-directed Lawton SN2' reaction
 establishes the required product, di-Me alc. II, stereochem. with high
 selectivity.

IT 906076-60-6 906076-63-9
 RL: RCT (Reactant); RACT (Reactant or reagent)
 (preparation of stereotetrads of aplyronine A via double Lawton nucleophilic
 substitution to epoxyvinyl sulfones)
RN 906076-60-6 CAPLUS
CN Silane, (1,1-dimethylethyl)dimethyl[[(1R,2S)-2-methyl-5-(phenylsulfonyl) 3,5-cycloheptadien-1-yl]oxy]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 906076-63-9 CAPLUS

CN Silane, (1,1-dimethylethyl)dimethyl[[(1S,2S)-2-methyl-5-(phenylsulfonyl)-3,5-cycloheptadien-1-yl]oxy]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 906076-69-5 CAPLUS

CN Silane, (1,1-dimethylethyl)dimethyl[[(1S,2R,3R,7S)-2-methyl-6-(phenylsulfonyl)-8-oxabicyclo[5.1.0]oct-5-en-3-yl]oxy]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 906076-70-8 CAPLUS

CN 2-Cyclohepten-1-ol, 6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4-(3,5-dimethyl-1H-pyrazol-1-yl)-7-methyl-3-(phenylsulfonyl)-, (1S,4R,6R,7S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 906076-72-0 CAPLUS

CN 3-Cyclohepten-1-ol, 6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-2,7-dimethyl-3-(phenylsulfonyl)-, (1S,2S,6R,7S)- (9CI) (CA INDEX NAME)

RN 906076-73-1 CAPLUS

CN 3-Cyclohepten-1-ol, 6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-2,7-dimethyl-3-(phenylsulfonyl)-, (1S,2S,6R,7R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 906076-78-6 CAPLUS

CN 3-Cyclohepten-1-ol, 6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-5,7-dimethyl-4-(phenylsulfonyl)-, (1R,5S,6S,7S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 906076-87-7 CAPLUS

CN 3-Cyclohepten-1-ol, 6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-2,7-dimethyl-3-(phenylsulfonyl)-, (1S,2R,6R,7S)- (9CI) (CA INDEX NAME)

IT 906076-91-3P

RL: SPN (Synthetic preparation); PREP (Preparation) (preparation of stereotetrads of aplyronine A via double Lawton nucleophilic substitution to epoxyvinyl sulfones)

RN 906076-91-3 CAPLUS

CN 3-Cyclohepten-1-ol, 6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-2,7-dimethyl-3-(phenylsulfonyl)-, (1S,2R,6R,7R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 906076-71-9P 906076-75-3P

RL: PRP (Properties); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of stereotetrads of aplyronine A via double Lawton nucleophilic substitution to epoxyvinyl sulfones and crystal structure)

RN 906076-71-9 CAPLUS

CN 2-Cyclohepten-1-ol, 6-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-4-(3,5-dimethyl-1H-pyrazol-1-yl)-7-methyl-3-(phenylsulfonyl)-, (1S,4R,6R,7R)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 906076-75-3 CAPLUS

CN Silane, [[(1R,2S,3S,4S)-2,4-dimethyl-5-(phenylsulfonyl)-5-cycloheptene-1,3-

Absolute stereochemistry.

IT 906076-74-2P 906076-90-2P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation) (preparation of stereotetrads of aplyronine A via double Lawton nucleophilic substitution to epoxyvinyl sulfones and crystal structure)

RN 906076-74-2 CAPLUS

CN Silane, [[(1R,2R,3S,4S)-2,4-dimethyl-5-(phenyIsulfonyl)-5-cycloheptene-1,3-diyl]bis(oxy)]bis[(1,1-dimethylethyl)dimethyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 906076-90-2 CAPLUS

CN Silane, [[(1R,2R,3S,4R)-2,4-dimethyl-5-(phenylsulfonyl)-5-cycloheptene-1,3-diyl]bis(oxy)]bis[(1,1-dimethylethyl)dimethyl- (9CI) (CA INDEX NAME)

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AUTHOR(S):

PUBLISHER:

145:166888

TITLE:

Asymmetric allylic alkylation of cyclic vinylogous esters and thioesters by Pd-catalyzed decarboxylation

of enol carbonate and β -keto ester substrates Trost, Barry M.; Bream, Robert N.; Xu, Jiayi Department of Chemistry, Stanford University,

CORPORATE SOURCE:

Stanford, CA, 94305-5080, USA

SOURCE:

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Excellent yields and enantioselectivities were achieved for the palladium-catalyzed asym. allylic alkylation of vinylogous thioesters. The close-to-neutral reaction conditions ensure that this reaction can tolerate a wide range of functionalities. Furthermore, this approach provides a convenient protocol for the synthesis of synthetically important α, α - and γ, γ -disubstituted cycloalkenones.

IT 900493-60-9

RL: RCT (Reactant); RACT (Reactant or reagent)

(asym. allylic alkylation of cyclic vinylogous esters and thioesters by Pd-catalyzed decarboxylation of enol carbonates and β -keto esters)

RN 900493-60-9 CAPLUS

Carbonic acid, 2-methyl-6-(phenylthio)-1,6-cycloheptadien-1-yl 2-propenyl CN ester (9CI) (CA INDEX NAME)

ΙT 900493-67-6P

RL: SPN (Synthetic preparation); PREP (Preparation)

(asym. allylic alkylation of cyclic vinylogous esters and thioesters by Pd-catalyzed decarboxylation of enol carbonates and β -keto esters)

900493-67-6 CAPLUS RN

CN 2-Cyclohepten-1-one, 7-methyl-3-(phenylthio)-7-(2-propenyl)- (9CI) (CA INDEX NAME)

PhS
$$CH_2-CH=CH_2$$

THERE ARE 27 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT